



Full wwPDB EM Validation Report ⓘ

Dec 1, 2021 – 04:03 pm GMT

PDB ID : 7OH3
EMDB ID : EMD-12892
Title : Nog1-TAP associated immature ribosomal particle population B from *S. cerevisiae*
Authors : Milkereit, P.; Poell, G.
Deposited on : 2021-05-08
Resolution : 3.40 Å(reported)
Based on initial model : 3JCT

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

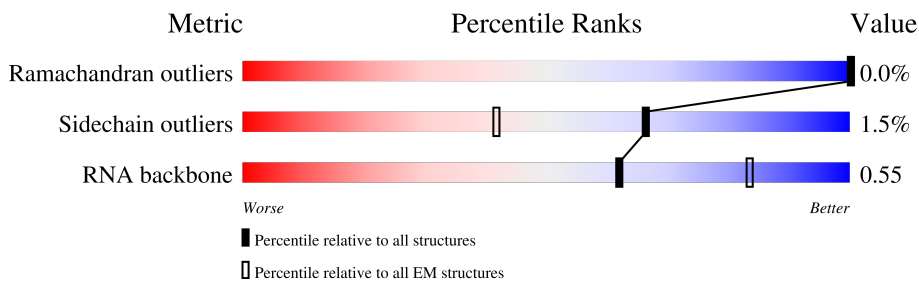
EMDB validation analysis : 0.0.0.dev97
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	1	3396	
2	2	158	
3	3	121	
4	5	120	
5	A	254	
6	B	387	
7	C	362	
8	D	297	

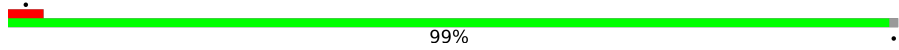

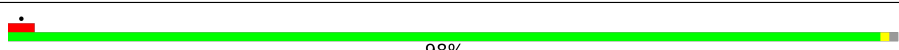
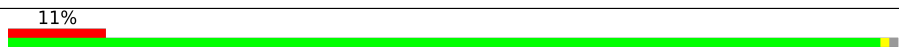
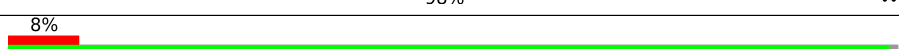

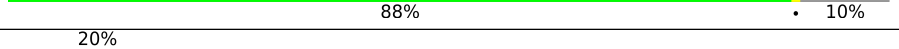
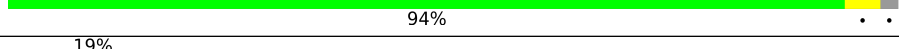
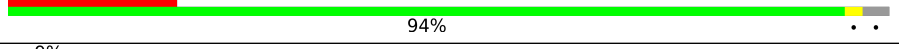
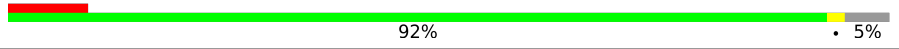


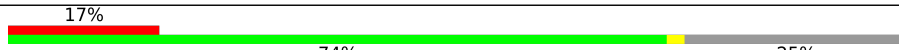


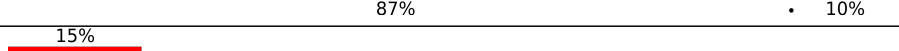
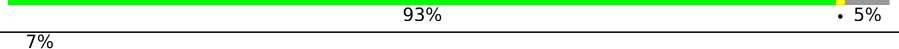
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Mol	Chain	Length	Quality of chain
9	E	176	89% 11%
10	F	244	91% 9%
11	G	256	11% 86% 11%
12	H	191	97% ..
13	J	174	11% 96% ..
14	L	199	9% 93% 6%
15	M	138	5% 98% ..
16	N	204	15% 99%
17	O	199	5% 98% ..
18	P	184	7% 93% 5%
19	Q	186	71% 28%
20	R	189	8% 81% 17%
21	S	172	10% 98% ..
22	T	160	12% 69% 29%
23	U	121	11% 80% 19%
24	V	137	8% 99% ..
25	W	236	97% ..
26	X	142	82% 18%
27	Y	127	98% ..
28	Z	136	9% 99% ..
29	a	149	62% 38%
30	b	647	6% 69% 30%
31	c	105	91% 8%
32	d	113	13% 93% 7%
33	e	130	15% 98%

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Mol	Chain	Length	Quality of chain
34	f	107	 99%
35	g	121	 92% 7%
36	h	120	 98%
37	i	100	 98%
38	j	88	 99%
39	k	78	 88% 10%
40	l	51	 94%
41	m	486	 94%
42	p	92	 92% 5%
43	r	261	 87% 12%
44	s	520	 90%
45	u	199	 74% 25%
46	v	344	 83% 17%
47	w	203	 87% 10%
48	x	515	 93% 5%
49	y	245	 98%
50	z	106	 40% 60%

2 Entry composition

There are 52 unique types of molecules in this entry. The entry contains 240945 atoms, of which 104342 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a RNA chain called 25S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
1	1	3013	96890	28804	32398	11667	21009	3012	0	0

- Molecule 2 is a RNA chain called 5.8S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
2	2	155	4957	1473	1664	580	1085	155	0	0

- Molecule 3 is a RNA chain called 5S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
3	3	121	3883	1152	1304	461	845	121	0	0

- Molecule 4 is a protein called rRNA-processing protein CGR1.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
4	5	73	1334	395	689	133	114	3	0	0

- Molecule 5 is a protein called 60S ribosomal protein L2-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
5	A	196	3098	954	1579	301	264	0	0

- Molecule 6 is a protein called 60S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
6	B	386	6247	1956	3166	584	533	8	0	0

- Molecule 7 is a protein called 60S ribosomal protein L4-A.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
7	C	352	Total	C	H	N	O	S	0	0
			5490	1691	2804	509	483	3		

- Molecule 8 is a protein called 60S ribosomal protein L5.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
8	D	267	Total	C	H	N	O	S	0	0
			4266	1361	2114	379	410	2		

- Molecule 9 is a protein called 60S ribosomal protein L6-A.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
9	E	156	Total	C	H	N	O	S	0	0
			2567	800	1328	222	216	1		

- Molecule 10 is a protein called 60S ribosomal protein L7-A.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
10	F	222	Total	C	H	N	O	S	0	0
			3647	1151	1863	324	308	1		

- Molecule 11 is a protein called 60S ribosomal protein L8-A.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
11	G	228	Total	C	H	N	O	S	0	0
			3668	1142	1884	320	319	3		

- Molecule 12 is a protein called 60S ribosomal protein L9-A.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
12	H	188	Total	C	H	N	O	S	0	0
			3059	948	1566	271	270	4		

- Molecule 13 is a protein called 60S ribosomal protein L11-A.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
13	J	169	Total	C	H	N	O	S	0	0
			2738	847	1385	253	249	4		

- Molecule 14 is a protein called 60S ribosomal protein L13-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
14	L	187	3057	934	1558	307	258	0	0

- Molecule 15 is a protein called 60S ribosomal protein L14-A.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
15	M	137	2214	678	1155	200	179	2	0	0

- Molecule 16 is a protein called 60S ribosomal protein L15-A.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
16	N	203	3500	1077	1780	361	281	1	0	0

- Molecule 17 is a protein called 60S ribosomal protein L16-A.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
17	O	197	3216	1003	1661	289	262	1	0	0

- Molecule 18 is a protein called 60S ribosomal protein L17-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
18	P	174	2798	857	1419	276	246	0	0

- Molecule 19 is a protein called 60S ribosomal protein L18-A.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
19	Q	134	2151	659	1116	196	179	1	0	0

- Molecule 20 is a protein called 60S ribosomal protein L19-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
20	R	156	2601	781	1343	265	212	0	0

- Molecule 21 is a protein called 60S ribosomal protein L20-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
21	S	171	2913	925	1476	266	243	3	0	0

- Molecule 22 is a protein called 60S ribosomal protein L21-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
22	T	113	1868	574	962	172	157	3	0	0

- Molecule 23 is a protein called 60S ribosomal protein L22-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
23	U	98	1571	502	796	128	145		0	0

- Molecule 24 is a protein called 60S ribosomal protein L23-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
24	V	136	2052	628	1049	189	179	7	0	0

- Molecule 25 is a protein called Ribosome assembly factor MRT4.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
25	W	234	3806	1194	1921	323	362	6	0	0

- Molecule 26 is a protein called 60S ribosomal protein L25.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
26	X	117	1932	602	995	164	169	2	0	0

- Molecule 27 is a protein called 60S ribosomal protein L26-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
27	Y	126	2075	625	1082	192	176		0	0

- Molecule 28 is a protein called 60S ribosomal protein L27-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
28	Z	135	2248	710	1156	202	180	0	0

- Molecule 29 is a protein called 60S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
29	a	93	1512	479	777	130	125	1	0	0

- Molecule 30 is a protein called Nucleolar GTP-binding protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
30	b	455	7450	2353	3755	640	684	18	0	0

- Molecule 31 is a protein called 60S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
31	c	97	1541	479	798	124	139	1	0	0

- Molecule 32 is a protein called 60S ribosomal protein L31-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
32	d	105	1761	544	905	163	148	1	0	0

- Molecule 33 is a protein called 60S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
33	e	127	2111	647	1091	205	167	1	0	0

- Molecule 34 is a protein called 60S ribosomal protein L33-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
34	f	106	1731	540	881	165	144	1	0	0

- Molecule 35 is a protein called 60S ribosomal protein L34-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
35	g	112	Total	C	H	N	O	S	0	0
			1831	546	950	179	152	4		

- Molecule 36 is a protein called 60S ribosomal protein L35-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
36	h	119	Total	C	H	N	O	S	0	0
			2048	615	1079	186	167	1		

- Molecule 37 is a protein called 60S ribosomal protein L36-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
37	i	99	Total	C	H	N	O	S	0	0
			1621	481	850	156	132	2		

- Molecule 38 is a protein called 60S ribosomal protein L37-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
38	j	87	Total	C	H	N	O	S	0	0
			1369	414	688	148	114	5		

- Molecule 39 is a protein called 60S ribosomal protein L38.

Mol	Chain	Residues	Atoms					AltConf	Trace	
39	k	70	Total	C	H	N	O	S	0	0
			1197	362	634	106	95			

- Molecule 40 is a protein called 60S ribosomal protein L39.

Mol	Chain	Residues	Atoms					AltConf	Trace	
40	l	50	Total	C	H	N	O	S	0	0
			912	272	476	97	65	2		

- Molecule 41 is a protein called Nucleolar GTP-binding protein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace	
41	m	469	Total	C	H	N	O	S	0	0
			7611	2381	3837	685	699	9		

- Molecule 42 is a protein called 60S ribosomal protein L43-A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
42	p	87	1375	411	710	134	115	5	0	0

- Molecule 43 is a protein called Ribosome biogenesis protein NSA2.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
43	r	230	3827	1177	1967	352	324	7	0	0

- Molecule 44 is a protein called Nuclear GTP-binding protein NUG1.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
44	s	54	947	279	503	90	73	2	0	0

- Molecule 45 is a protein called Ribosome biogenesis protein RLP24.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
45	u	150	2582	793	1317	253	210	9	0	0

- Molecule 46 is a protein called Ribosome biogenesis protein RPF2.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
46	v	287	4718	1482	2400	408	412	16	0	0

- Molecule 47 is a protein called Regulator of ribosome biosynthesis.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
47	w	182	2960	911	1512	261	271	5	0	0

- Molecule 48 is a protein called Ribosome assembly protein 4.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
48	x	488	7606	2398	3799	677	711	21	0	0

- Molecule 49 is a protein called Eukaryotic translation initiation factor 6.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
49	y	244	3685	1146	1836	319	377	7	0	0

- Molecule 50 is a protein called UPF0642 protein YBL028C.

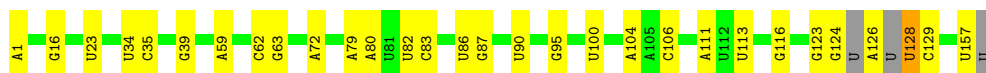
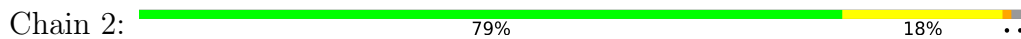
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
50	z	42	699	206	364	69	60	0	0

- Molecule 51 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

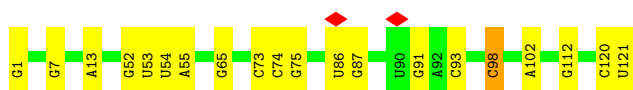
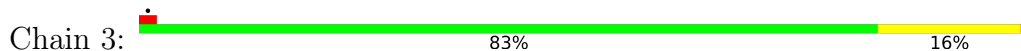
Mol	Chain	Residues	Atoms		AltConf
			Total	Mg	
51	b	1	1	1	0
51	m	1	1	1	0

- Molecule 52 is ZINC ION (three-letter code: ZN) (formula: Zn).

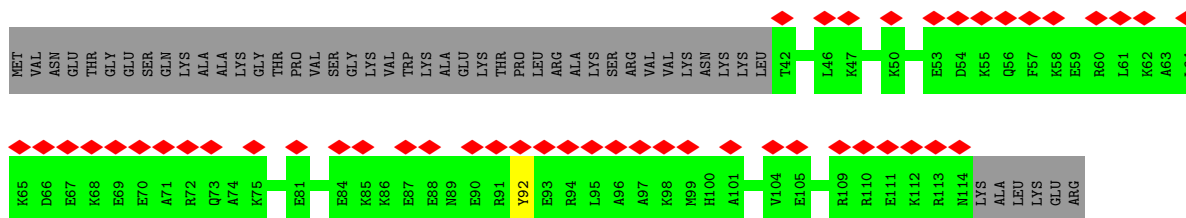
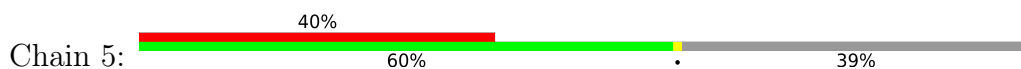
Mol	Chain	Residues	Atoms		AltConf
			Total	Zn	
52	j	1	1	1	0
52	p	1	1	1	0
52	u	1	1	1	0



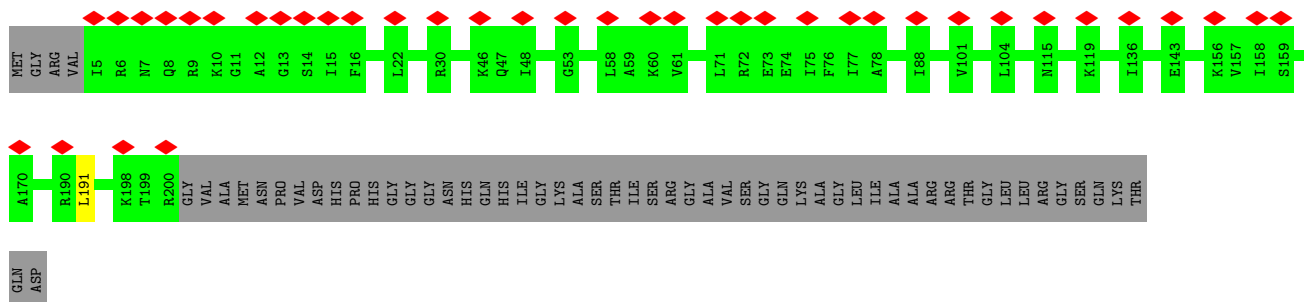
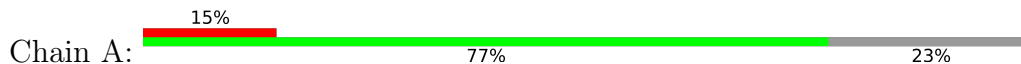
• Molecule 3: 5S rRNA



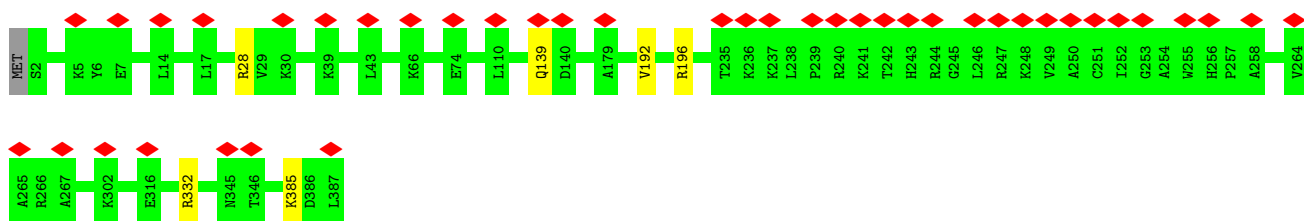
• Molecule 4: rRNA-processing protein CGR1



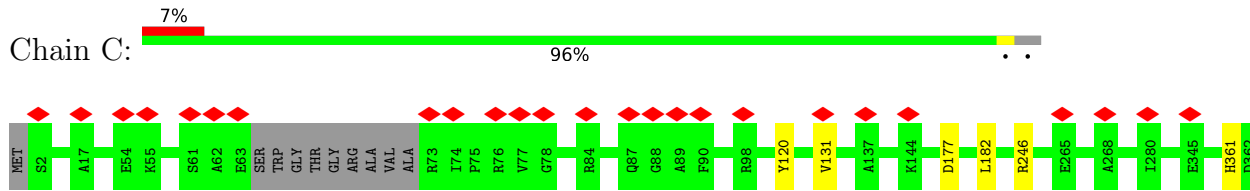
• Molecule 5: 60S ribosomal protein L2-A



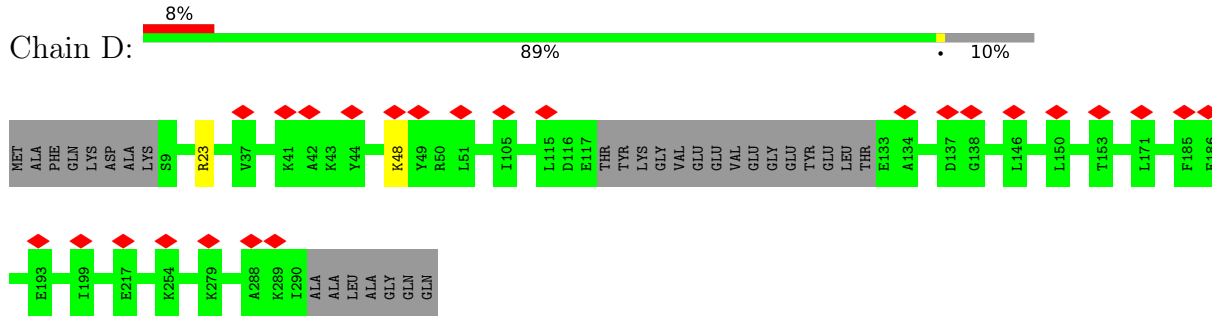
• Molecule 6: 60S ribosomal protein L3



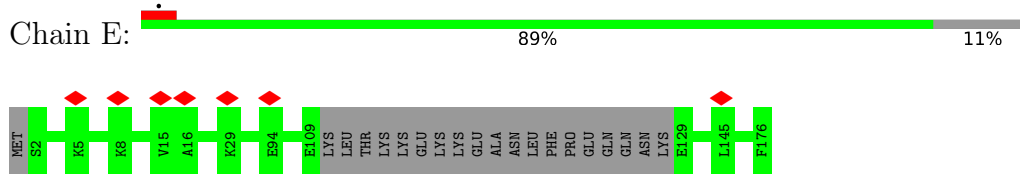
• Molecule 7: 60S ribosomal protein L4-A



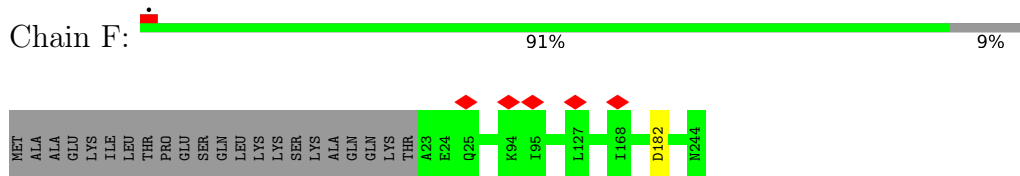
- Molecule 8: 60S ribosomal protein L5



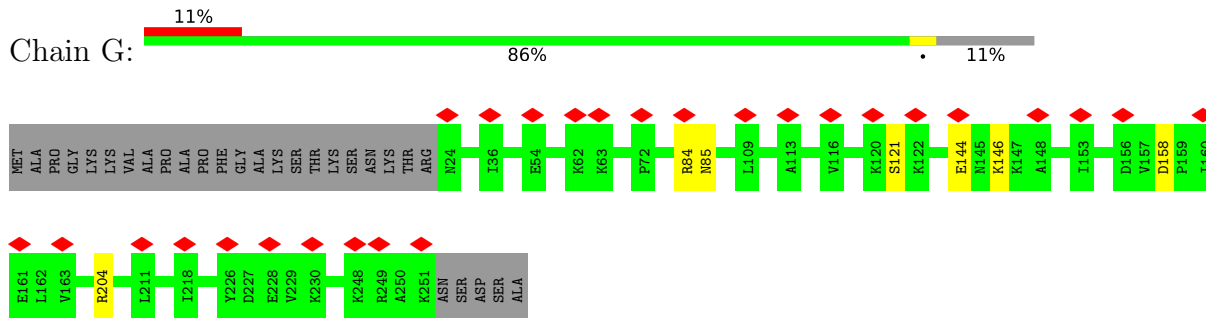
- Molecule 9: 60S ribosomal protein L6-A



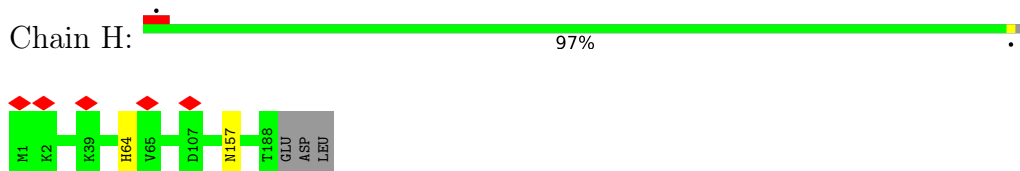
- Molecule 10: 60S ribosomal protein L7-A



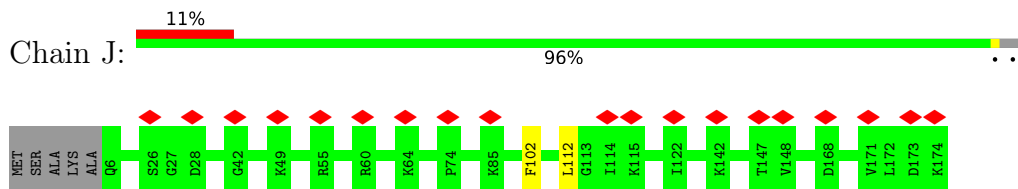
- Molecule 11: 60S ribosomal protein L8-A



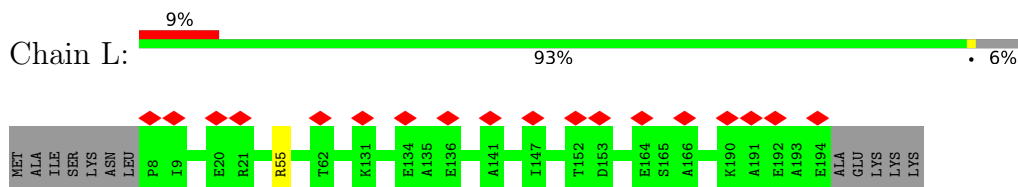
- Molecule 12: 60S ribosomal protein L9-A



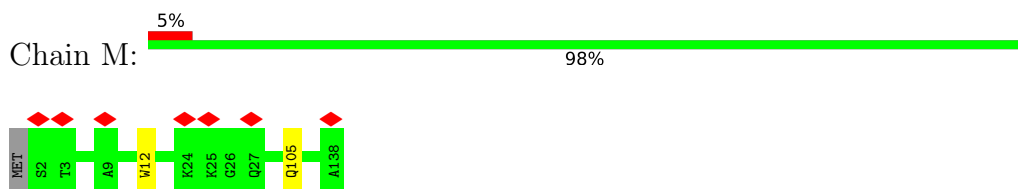
- Molecule 13: 60S ribosomal protein L11-A



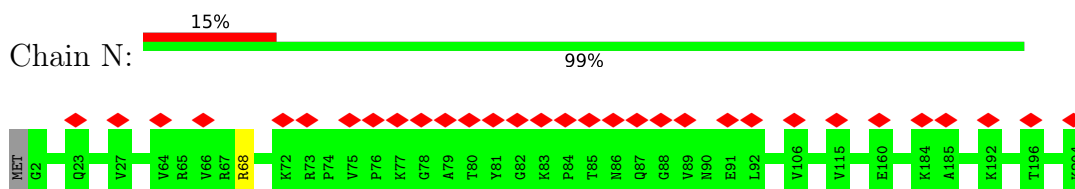
- Molecule 14: 60S ribosomal protein L13-A



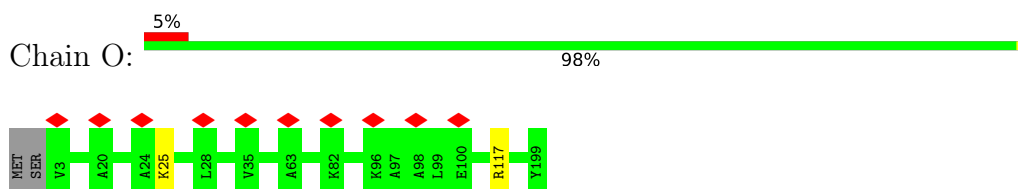
- Molecule 15: 60S ribosomal protein L14-A



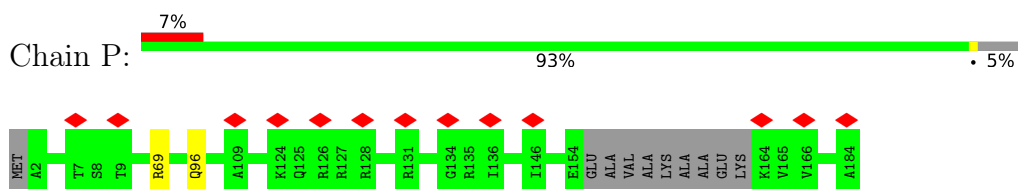
- Molecule 16: 60S ribosomal protein L15-A



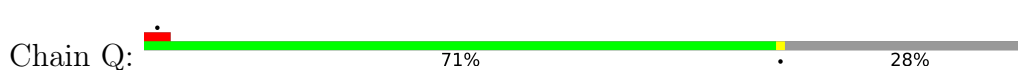
- Molecule 17: 60S ribosomal protein L16-A

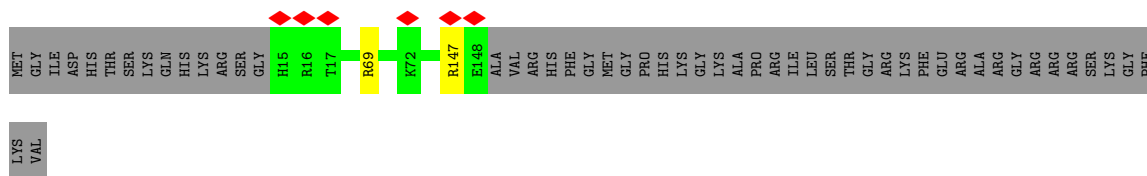


- Molecule 18: 60S ribosomal protein L17-A

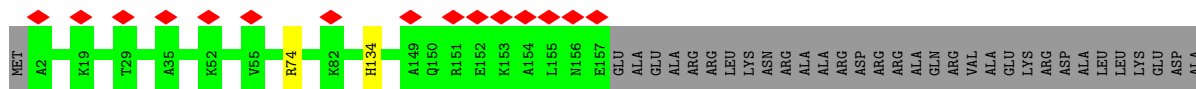
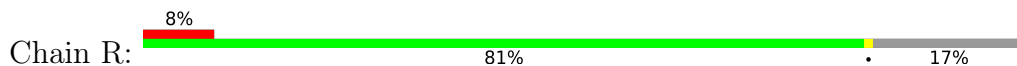


- Molecule 19: 60S ribosomal protein L18-A

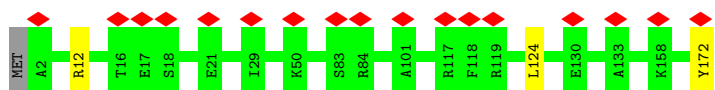




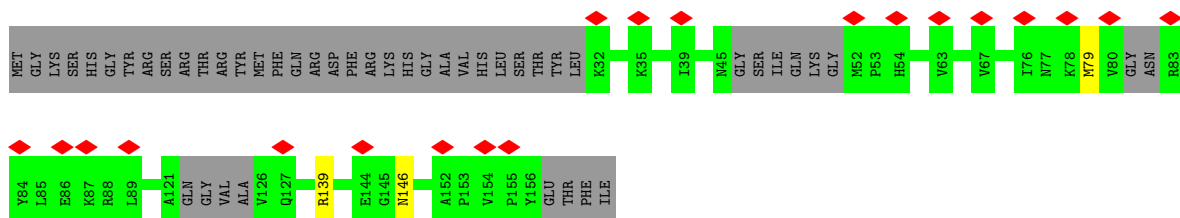
• Molecule 20: 60S ribosomal protein L19-A



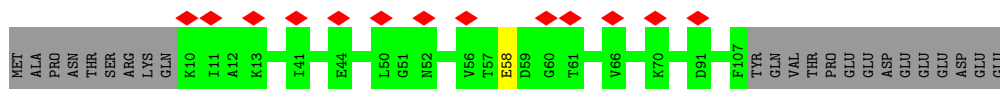
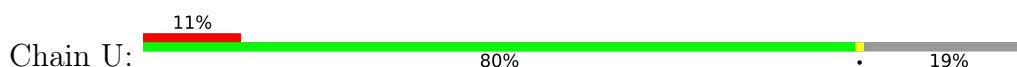
• Molecule 21: 60S ribosomal protein L20-A



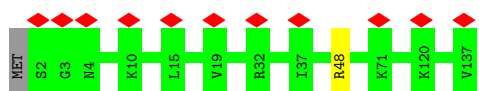
• Molecule 22: 60S ribosomal protein L21-A



• Molecule 23: 60S ribosomal protein L22-A

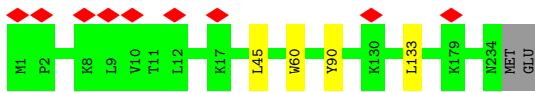


• Molecule 24: 60S ribosomal protein L23-A




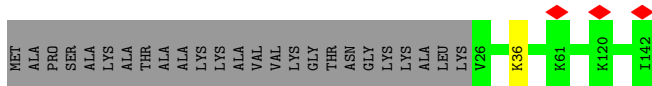
• Molecule 25: Ribosome assembly factor MRT4

Chain W:  97%



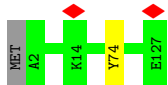
- Molecule 26: 60S ribosomal protein L25

Chain X:  82%



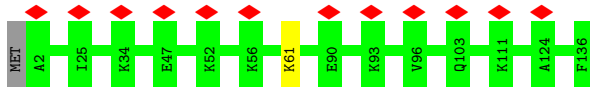
- Molecule 27: 60S ribosomal protein L26-A

Chain Y:  98%



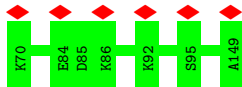
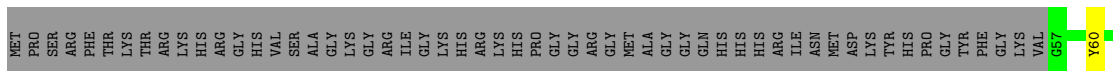
- Molecule 28: 60S ribosomal protein L27-A

Chain Z:  99%



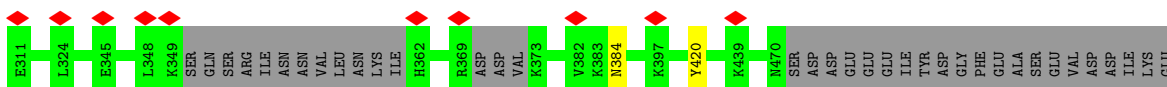
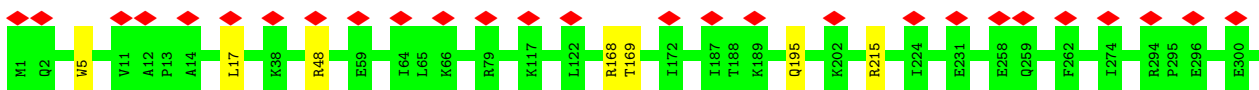
- Molecule 29: 60S ribosomal protein L28

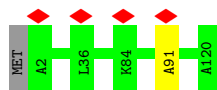
Chain a:  62%



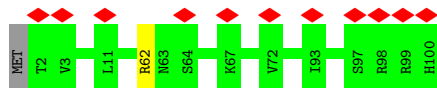
- Molecule 30: Nucleolar GTP-binding protein 1

Chain b:  69%

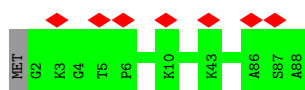




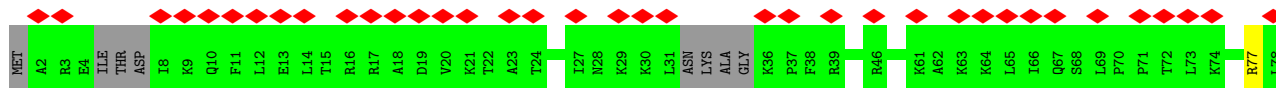
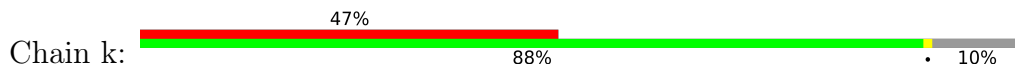
- Molecule 37: 60S ribosomal protein L36-A



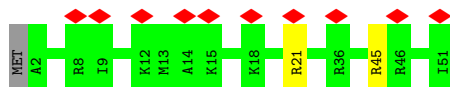
- Molecule 38: 60S ribosomal protein L37-A



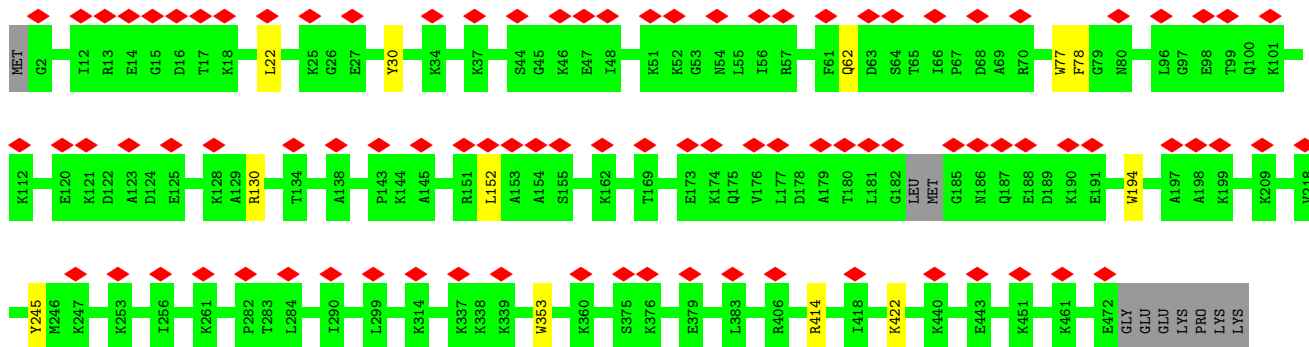
- Molecule 39: 60S ribosomal protein L38

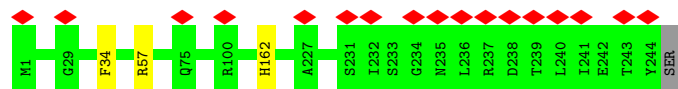


- Molecule 40: 60S ribosomal protein L39

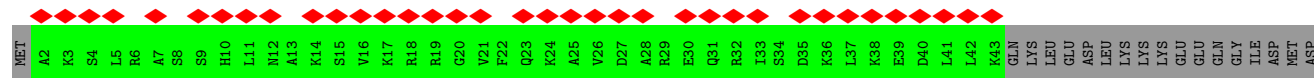


- Molecule 41: Nucleolar GTP-binding protein 2





• Molecule 50: UPF0642 protein YBL028C



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	18823	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	84.67	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	0.136	Depositor
Minimum map value	-0.051	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.008	Depositor
Recommended contour level	0.026	Depositor
Map size (Å)	425.40002, 425.40002, 425.40002	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.0635, 1.0635, 1.0635	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	1	0.16	0/72171	0.74	44/112459 (0.0%)
2	2	0.16	0/3678	0.77	4/5722 (0.1%)
3	3	0.16	0/2883	0.78	6/4491 (0.1%)
4	5	0.24	0/649	0.37	0/848
5	A	0.24	0/1547	0.43	0/2079
6	B	0.24	0/3152	0.43	0/4239
7	C	0.24	0/2735	0.41	0/3700
8	D	0.24	0/2198	0.40	0/2964
9	E	0.25	0/1260	0.40	0/1694
10	F	0.25	0/1821	0.38	0/2451
11	G	0.24	0/1816	0.41	0/2450
12	H	0.24	0/1514	0.42	0/2039
13	J	0.24	0/1374	0.43	0/1842
14	L	0.24	0/1524	0.41	0/2046
15	M	0.23	0/1074	0.40	0/1446
16	N	0.23	0/1757	0.41	0/2354
17	O	0.24	0/1585	0.38	0/2128
18	P	0.24	0/1401	0.41	0/1881
19	Q	0.25	0/1050	0.42	0/1419
20	R	0.23	0/1275	0.39	0/1702
21	S	0.24	0/1473	0.41	0/1980
22	T	0.24	0/918	0.41	0/1229
23	U	0.25	0/790	0.43	0/1069
24	V	0.25	0/1018	0.42	0/1369
25	W	0.24	0/1918	0.41	0/2586
26	X	0.24	0/952	0.40	0/1285
27	Y	0.23	0/1004	0.39	0/1341
28	Z	0.24	0/1118	0.41	0/1497
29	a	0.25	0/751	0.39	0/1013
30	b	0.24	0/3764	0.40	0/5075
31	c	0.24	0/751	0.39	0/1008
32	d	0.23	0/870	0.39	0/1168

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	e	0.23	0/1041	0.40	0/1394
34	f	0.25	0/868	0.42	0/1168
35	g	0.23	0/891	0.42	0/1191
36	h	0.24	0/978	0.39	0/1301
37	i	0.24	0/778	0.38	0/1034
38	j	0.25	0/696	0.42	0/923
39	k	0.25	0/567	0.42	0/754
40	l	0.24	0/443	0.41	0/588
41	m	0.23	0/3848	0.41	0/5181
42	p	0.24	0/672	0.45	0/895
43	r	0.24	0/1892	0.42	0/2528
44	s	0.24	0/448	0.39	0/585
45	u	0.24	0/1287	0.38	0/1711
46	v	0.24	0/2361	0.40	0/3153
47	w	0.23	0/1471	0.40	0/1980
48	x	0.23	0/3897	0.41	0/5282
49	y	0.23	0/1872	0.43	0/2548
50	z	0.24	0/336	0.33	0/443
All	All	0.20	0/146137	0.62	54/213233 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
41	m	0	1

There are no bond length outliers.

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	3	98	C	C2-N1-C1'	7.20	126.72	118.80
1	1	1082	U	OP1-P-OP2	-6.81	109.39	119.60
1	1	2376	G	OP1-P-OP2	-6.80	109.39	119.60
1	1	3157	U	OP1-P-OP2	-6.80	109.40	119.60
1	1	2943	G	OP1-P-OP2	-6.80	109.41	119.60
1	1	2645	G	OP1-P-OP2	-6.79	109.41	119.60
2	2	128	U	OP1-P-OP2	-6.79	109.42	119.60
1	1	2313	A	OP1-P-OP2	-6.79	109.42	119.60
1	1	1353	U	OP1-P-OP2	-6.78	109.43	119.60
3	3	1	G	OP1-P-OP2	-6.78	109.43	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	547	G	OP1-P-OP2	-6.78	109.43	119.60
1	1	2540	A	OP1-P-OP2	-6.78	109.44	119.60
2	2	1	A	OP1-P-OP2	-6.78	109.44	119.60
1	1	2765	C	OP1-P-OP2	-6.77	109.44	119.60
1	1	1	G	OP1-P-OP2	-6.77	109.45	119.60
1	1	1047	A	OP1-P-OP2	-6.76	109.45	119.60
1	1	2619	G	OP1-P-OP2	-6.76	109.45	119.60
1	1	1200	A	OP1-P-OP2	-6.76	109.46	119.60
1	1	2571	U	OP1-P-OP2	-6.76	109.46	119.60
2	2	126	A	OP1-P-OP2	-6.76	109.46	119.60
1	1	1575	A	OP1-P-OP2	-6.76	109.47	119.60
1	1	2979	U	OP1-P-OP2	-6.76	109.47	119.60
1	1	2501	U	OP1-P-OP2	-6.75	109.47	119.60
1	1	2946	A	OP1-P-OP2	-6.75	109.48	119.60
1	1	3	U	OP1-P-OP2	-6.75	109.48	119.60
1	1	253	A	OP1-P-OP2	-6.75	109.48	119.60
1	1	2194	G	OP1-P-OP2	-6.74	109.48	119.60
1	1	1096	U	OP1-P-OP2	-6.74	109.49	119.60
1	1	2093	A	OP1-P-OP2	-6.74	109.49	119.60
1	1	2860	U	OP1-P-OP2	-6.73	109.50	119.60
1	1	2847	A	OP1-P-OP2	-6.72	109.51	119.60
1	1	2411	U	OP1-P-OP2	-6.72	109.52	119.60
1	1	2802	A	OP1-P-OP2	-6.72	109.52	119.60
1	1	494	G	OP1-P-OP2	-6.72	109.52	119.60
1	1	2260	U	OP1-P-OP2	-6.72	109.53	119.60
1	1	3356	G	OP1-P-OP2	-6.71	109.53	119.60
1	1	1765	U	OP1-P-OP2	-6.71	109.54	119.60
1	1	2586	G	OP1-P-OP2	-6.70	109.56	119.60
3	3	73	C	C2-N1-C1'	6.67	126.13	118.80
1	1	1283	C	N3-C2-O2	-6.25	117.53	121.90
1	1	2654	C	C2-N1-C1'	6.12	125.53	118.80
1	1	2798	C	N3-C2-O2	-6.06	117.66	121.90
1	1	2979	U	N1-C1'-C2'	5.98	121.77	114.00
3	3	98	C	N1-C2-O2	5.97	122.48	118.90
1	1	2552	C	C2-N1-C1'	5.90	125.29	118.80
1	1	1283	C	N1-C2-O2	5.88	122.43	118.90
2	2	157	U	C2-N1-C1'	5.87	124.74	117.70
1	1	922	U	C2-N1-C1'	5.76	124.61	117.70
1	1	2550	U	C2-N1-C1'	5.75	124.61	117.70
1	1	2861	U	C2-N1-C1'	5.74	124.58	117.70
3	3	73	C	N1-C2-O2	5.71	122.33	118.90
1	1	3058	U	C2-N1-C1'	5.61	124.43	117.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	2112	U	C2-N1-C1'	5.59	124.41	117.70
3	3	98	C	N3-C2-O2	-5.29	118.20	121.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
41	m	77	TRP	Peptide

5.2 Too-close contacts [i](#)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	5	71/120 (59%)	71 (100%)	0	0	100	100
5	A	194/254 (76%)	187 (96%)	7 (4%)	0	100	100
6	B	384/387 (99%)	362 (94%)	22 (6%)	0	100	100
7	C	348/362 (96%)	323 (93%)	24 (7%)	1 (0%)	41	72
8	D	263/297 (89%)	248 (94%)	15 (6%)	0	100	100
9	E	152/176 (86%)	145 (95%)	7 (5%)	0	100	100
10	F	220/244 (90%)	215 (98%)	5 (2%)	0	100	100
11	G	226/256 (88%)	213 (94%)	13 (6%)	0	100	100
12	H	186/191 (97%)	178 (96%)	8 (4%)	0	100	100
13	J	167/174 (96%)	155 (93%)	12 (7%)	0	100	100
14	L	185/199 (93%)	169 (91%)	16 (9%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
15	M	135/138 (98%)	131 (97%)	4 (3%)	0	100	100
16	N	201/204 (98%)	193 (96%)	8 (4%)	0	100	100
17	O	195/199 (98%)	195 (100%)	0	0	100	100
18	P	170/184 (92%)	166 (98%)	4 (2%)	0	100	100
19	Q	132/186 (71%)	130 (98%)	2 (2%)	0	100	100
20	R	154/189 (82%)	153 (99%)	1 (1%)	0	100	100
21	S	169/172 (98%)	160 (95%)	9 (5%)	0	100	100
22	T	105/160 (66%)	96 (91%)	9 (9%)	0	100	100
23	U	96/121 (79%)	95 (99%)	1 (1%)	0	100	100
24	V	134/137 (98%)	133 (99%)	1 (1%)	0	100	100
25	W	232/236 (98%)	223 (96%)	9 (4%)	0	100	100
26	X	115/142 (81%)	112 (97%)	3 (3%)	0	100	100
27	Y	124/127 (98%)	122 (98%)	2 (2%)	0	100	100
28	Z	133/136 (98%)	125 (94%)	8 (6%)	0	100	100
29	a	91/149 (61%)	89 (98%)	2 (2%)	0	100	100
30	b	449/647 (69%)	421 (94%)	28 (6%)	0	100	100
31	c	95/105 (90%)	94 (99%)	1 (1%)	0	100	100
32	d	103/113 (91%)	100 (97%)	3 (3%)	0	100	100
33	e	125/130 (96%)	122 (98%)	3 (2%)	0	100	100
34	f	104/107 (97%)	98 (94%)	6 (6%)	0	100	100
35	g	110/121 (91%)	109 (99%)	1 (1%)	0	100	100
36	h	117/120 (98%)	109 (93%)	7 (6%)	1 (1%)	17	49
37	i	97/100 (97%)	91 (94%)	6 (6%)	0	100	100
38	j	85/88 (97%)	84 (99%)	1 (1%)	0	100	100
39	k	64/78 (82%)	64 (100%)	0	0	100	100
40	l	48/51 (94%)	48 (100%)	0	0	100	100
41	m	465/486 (96%)	429 (92%)	35 (8%)	1 (0%)	47	78
42	p	85/92 (92%)	80 (94%)	5 (6%)	0	100	100
43	r	224/261 (86%)	206 (92%)	18 (8%)	0	100	100
44	s	50/520 (10%)	47 (94%)	3 (6%)	0	100	100
45	u	148/199 (74%)	143 (97%)	5 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
46	v	283/344 (82%)	277 (98%)	6 (2%)	0	100	100
47	w	178/203 (88%)	168 (94%)	10 (6%)	0	100	100
48	x	476/515 (92%)	456 (96%)	20 (4%)	0	100	100
49	y	242/245 (99%)	236 (98%)	6 (2%)	0	100	100
50	z	40/106 (38%)	38 (95%)	2 (5%)	0	100	100
All	All	8170/9771 (84%)	7809 (96%)	358 (4%)	3 (0%)	100	100

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
36	h	91	ALA
41	m	78	PHE
7	C	131	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	5	67/106 (63%)	66 (98%)	1 (2%)	65	82
5	A	155/196 (79%)	154 (99%)	1 (1%)	86	94
6	B	322/323 (100%)	316 (98%)	6 (2%)	57	78
7	C	283/289 (98%)	278 (98%)	5 (2%)	59	79
8	D	223/245 (91%)	221 (99%)	2 (1%)	78	90
9	E	134/153 (88%)	134 (100%)	0	100	100
10	F	186/205 (91%)	185 (100%)	1 (0%)	88	94
11	G	187/208 (90%)	180 (96%)	7 (4%)	34	62
12	H	168/171 (98%)	166 (99%)	2 (1%)	71	85
13	J	147/150 (98%)	145 (99%)	2 (1%)	67	83
14	L	149/159 (94%)	148 (99%)	1 (1%)	84	92
15	M	108/109 (99%)	106 (98%)	2 (2%)	57	78

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
16	N	175/176 (99%)	174 (99%)	1 (1%)	86	94
17	O	160/162 (99%)	158 (99%)	2 (1%)	69	84
18	P	140/146 (96%)	138 (99%)	2 (1%)	67	83
19	Q	110/151 (73%)	108 (98%)	2 (2%)	59	79
20	R	129/154 (84%)	127 (98%)	2 (2%)	62	81
21	S	155/156 (99%)	152 (98%)	3 (2%)	57	78
22	T	99/137 (72%)	96 (97%)	3 (3%)	41	68
23	U	85/107 (79%)	84 (99%)	1 (1%)	71	85
24	V	104/105 (99%)	103 (99%)	1 (1%)	76	88
25	W	211/213 (99%)	207 (98%)	4 (2%)	57	78
26	X	102/118 (86%)	101 (99%)	1 (1%)	76	88
27	Y	109/110 (99%)	108 (99%)	1 (1%)	78	90
28	Z	115/116 (99%)	114 (99%)	1 (1%)	78	90
29	a	76/119 (64%)	75 (99%)	1 (1%)	69	84
30	b	409/573 (71%)	400 (98%)	9 (2%)	52	75
31	c	81/88 (92%)	80 (99%)	1 (1%)	71	85
32	d	92/97 (95%)	92 (100%)	0	100	100
33	e	109/111 (98%)	109 (100%)	0	100	100
34	f	90/91 (99%)	90 (100%)	0	100	100
35	g	95/103 (92%)	94 (99%)	1 (1%)	73	86
36	h	104/105 (99%)	104 (100%)	0	100	100
37	i	81/82 (99%)	80 (99%)	1 (1%)	71	85
38	j	70/71 (99%)	70 (100%)	0	100	100
39	k	63/69 (91%)	62 (98%)	1 (2%)	62	81
40	l	45/46 (98%)	43 (96%)	2 (4%)	28	58
41	m	413/428 (96%)	403 (98%)	10 (2%)	49	74
42	p	68/72 (94%)	66 (97%)	2 (3%)	42	69
43	r	203/229 (89%)	199 (98%)	4 (2%)	55	77
44	s	48/445 (11%)	45 (94%)	3 (6%)	18	47
45	u	133/180 (74%)	130 (98%)	3 (2%)	50	74
46	v	258/309 (84%)	256 (99%)	2 (1%)	81	91

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
47	w	161/179 (90%)	156 (97%)	5 (3%)	40	68
48	x	428/451 (95%)	421 (98%)	7 (2%)	62	81
49	y	210/211 (100%)	207 (99%)	3 (1%)	67	83
50	z	36/95 (38%)	36 (100%)	0	100	100
All	All	7096/8319 (85%)	6987 (98%)	109 (2%)	66	82

All (109) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	5	92	TYR
5	A	191	LEU
6	B	28	ARG
6	B	139	GLN
6	B	192	VAL
6	B	196	ARG
6	B	332	ARG
6	B	385	LYS
7	C	120	TYR
7	C	177	ASP
7	C	182	LEU
7	C	246	ARG
7	C	361	HIS
8	D	23	ARG
8	D	48	LYS
10	F	182	ASP
11	G	84	ARG
11	G	85	ASN
11	G	121	SER
11	G	144	GLU
11	G	146	LYS
11	G	158	ASP
11	G	204	ARG
12	H	64	HIS
12	H	157	ASN
13	J	102	PHE
13	J	112	LEU
14	L	55	ARG
15	M	12	TRP
15	M	105	GLN
16	N	68	ARG
17	O	25	LYS

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Mol	Chain	Res	Type
17	O	117	ARG
18	P	69	ARG
18	P	96	GLN
19	Q	69	ARG
19	Q	147	ARG
20	R	74	ARG
20	R	134	HIS
21	S	12	ARG
21	S	124	LEU
21	S	172	TYR
22	T	79	MET
22	T	139	ARG
22	T	146	ASN
23	U	58	GLU
24	V	48	ARG
25	W	45	LEU
25	W	60	TRP
25	W	90	TYR
25	W	133	LEU
26	X	36	LYS
27	Y	74	TYR
28	Z	61	LYS
29	a	60	TYR
30	b	5	TRP
30	b	17	LEU
30	b	48	ARG
30	b	168	ARG
30	b	169	THR
30	b	195	GLN
30	b	215	ARG
30	b	384	ASN
30	b	420	TYR
31	c	83	LYS
35	g	102	LYS
37	i	62	ARG
39	k	77	ARG
40	l	21	ARG
40	l	45	ARG
41	m	22	LEU
41	m	30	TYR
41	m	62	GLN
41	m	130	ARG

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Mol	Chain	Res	Type
41	m	152	LEU
41	m	194	TRP
41	m	245	TYR
41	m	353	TRP
41	m	414	ARG
41	m	422	LYS
42	p	49	ARG
42	p	84	ARG
43	r	23	ARG
43	r	186	HIS
43	r	224	ASN
43	r	251	ARG
44	s	1	MET
44	s	2	ARG
44	s	33	MET
45	u	2	ARG
45	u	80	ARG
45	u	82	ASN
46	v	11	ARG
46	v	62	ARG
47	w	70	MET
47	w	83	GLN
47	w	111	MET
47	w	145	TRP
47	w	162	VAL
48	x	20	ARG
48	x	293	TYR
48	x	322	TRP
48	x	360	TYR
48	x	404	LEU
48	x	406	ASN
48	x	410	PHE
49	y	34	PHE
49	y	57	ARG
49	y	162	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
6	B	198	HIS
7	C	48	GLN
7	C	114	ASN

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Mol	Chain	Res	Type
7	C	221	ASN
7	C	260	GLN
11	G	38	GLN
11	G	59	GLN
11	G	85	ASN
11	G	138	HIS
11	G	145	ASN
11	G	240	ASN
13	J	95	ASN
13	J	150	ASN
17	O	50	ASN
17	O	122	GLN
18	P	96	GLN
21	S	157	GLN
23	U	52	ASN
23	U	87	ASN
28	Z	36	HIS
30	b	70	ASN
30	b	152	GLN
32	d	57	GLN
32	d	80	ASN
32	d	87	ASN
41	m	227	HIS
43	r	10	HIS
43	r	224	ASN
45	u	82	ASN
46	v	9	ASN
46	v	64	ASN
47	w	195	GLN
48	x	116	GLN
48	x	208	ASN
48	x	321	HIS
49	y	82	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	1	2987/3396 (87%)	579 (19%)	80 (2%)
2	2	153/158 (96%)	25 (16%)	2 (1%)
3	3	120/121 (99%)	17 (14%)	1 (0%)
All	All	3260/3675 (88%)	621 (19%)	83 (2%)

All (621) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	1	13	A
1	1	14	U
1	1	20	A
1	1	26	A
1	1	40	A
1	1	41	G
1	1	43	A
1	1	49	A
1	1	60	A
1	1	65	A
1	1	66	A
1	1	74	G
1	1	76	G
1	1	92	G
1	1	93	C
1	1	96	G
1	1	110	G
1	1	111	C
1	1	116	A
1	1	122	A
1	1	135	C
1	1	136	G
1	1	148	G
1	1	153	U
1	1	156	G
1	1	157	A
1	1	161	G
1	1	172	G
1	1	187	A
1	1	190	U
1	1	191	U
1	1	200	C
1	1	206	G
1	1	207	U
1	1	210	U
1	1	213	A
1	1	218	G
1	1	219	A
1	1	221	A
1	1	222	A
1	1	234	G
1	1	240	U

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Mol	Chain	Res	Type
1	1	241	G
1	1	243	G
1	1	249	U
1	1	250	U
1	1	251	G
1	1	266	A
1	1	269	G
1	1	284	A
1	1	285	A
1	1	286	U
1	1	305	U
1	1	315	C
1	1	323	A
1	1	325	A
1	1	329	U
1	1	339	C
1	1	342	A
1	1	352	A
1	1	375	A
1	1	376	G
1	1	398	A
1	1	401	U
1	1	402	A
1	1	403	C
1	1	404	G
1	1	421	G
1	1	422	A
1	1	429	U
1	1	438	A
1	1	439	C
1	1	440	A
1	1	521	A
1	1	535	G
1	1	548	G
1	1	552	G
1	1	557	A
1	1	559	A
1	1	578	A
1	1	579	G
1	1	589	A
1	1	592	A
1	1	604	G

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Mol	Chain	Res	Type
1	1	611	A
1	1	619	A
1	1	620	U
1	1	621	A
1	1	622	A
1	1	643	U
1	1	644	G
1	1	645	A
1	1	646	A
1	1	647	A
1	1	648	C
1	1	650	C
1	1	677	A
1	1	681	U
1	1	691	A
1	1	699	A
1	1	705	A
1	1	716	A
1	1	720	A
1	1	721	G
1	1	722	G
1	1	735	A
1	1	742	G
1	1	761	A
1	1	764	U
1	1	765	C
1	1	766	U
1	1	767	U
1	1	768	C
1	1	776	U
1	1	777	U
1	1	780	A
1	1	781	G
1	1	785	G
1	1	787	G
1	1	806	A
1	1	817	A
1	1	830	A
1	1	836	A
1	1	850	U
1	1	857	G
1	1	861	C

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Mol	Chain	Res	Type
1	1	875	G
1	1	879	U
1	1	880	G
1	1	896	A
1	1	907	G
1	1	908	G
1	1	909	G
1	1	914	A
1	1	916	G
1	1	917	A
1	1	921	A
1	1	944	C
1	1	953	G
1	1	959	C
1	1	964	G
1	1	965	A
1	1	970	A
1	1	974	G
1	1	975	C
1	1	976	U
1	1	977	C
1	1	978	G
1	1	979	U
1	1	980	A
1	1	981	U
1	1	991	G
1	1	992	A
1	1	993	G
1	1	994	G
1	1	995	U
1	1	997	A
1	1	998	A
1	1	999	G
1	1	1000	C
1	1	1003	A
1	1	1048	A
1	1	1050	U
1	1	1056	U
1	1	1058	U
1	1	1063	G
1	1	1064	A
1	1	1065	A

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Mol	Chain	Res	Type
1	1	1097	G
1	1	1098	A
1	1	1103	A
1	1	1104	G
1	1	1105	A
1	1	1111	U
1	1	1115	G
1	1	1116	G
1	1	1117	G
1	1	1118	C
1	1	1123	U
1	1	1129	A
1	1	1132	C
1	1	1153	A
1	1	1155	C
1	1	1180	A
1	1	1181	U
1	1	1191	U
1	1	1192	C
1	1	1193	A
1	1	1196	C
1	1	1197	A
1	1	1204	A
1	1	1206	G
1	1	1207	G
1	1	1208	U
1	1	1209	G
1	1	1220	U
1	1	1221	A
1	1	1222	G
1	1	1223	A
1	1	1234	G
1	1	1240	A
1	1	1241	U
1	1	1242	G
1	1	1245	A
1	1	1251	A
1	1	1252	A
1	1	1254	C
1	1	1257	C
1	1	1259	A
1	1	1263	A

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Mol	Chain	Res	Type
1	1	1266	G
1	1	1271	A
1	1	1272	C
1	1	1278	A
1	1	1286	A
1	1	1287	A
1	1	1302	A
1	1	1303	A
1	1	1304	A
1	1	1307	G
1	1	1330	A
1	1	1348	U
1	1	1354	G
1	1	1356	U
1	1	1357	G
1	1	1386	A
1	1	1399	A
1	1	1400	G
1	1	1406	A
1	1	1408	G
1	1	1417	G
1	1	1419	A
1	1	1434	G
1	1	1436	U
1	1	1437	C
1	1	1455	U
1	1	1470	U
1	1	1481	A
1	1	1482	A
1	1	1496	C
1	1	1508	C
1	1	1539	A
1	1	1555	U
1	1	1557	A
1	1	1561	G
1	1	1562	C
1	1	1563	C
1	1	1581	C
1	1	1582	C
1	1	1583	A
1	1	1589	A
1	1	1593	A

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Mol	Chain	Res	Type
1	1	1607	U
1	1	1608	C
1	1	1620	U
1	1	1629	U
1	1	1639	C
1	1	1642	A
1	1	1643	A
1	1	1645	U
1	1	1647	A
1	1	1657	C
1	1	1683	A
1	1	1688	U
1	1	1705	U
1	1	1713	G
1	1	1717	U
1	1	1724	U
1	1	1725	C
1	1	1742	U
1	1	1750	A
1	1	1751	G
1	1	1766	G
1	1	1770	G
1	1	1796	G
1	1	1797	A
1	1	1807	G
1	1	1808	G
1	1	1812	G
1	1	1814	A
1	1	1816	A
1	1	1817	G
1	1	1820	U
1	1	1821	U
1	1	1839	A
1	1	1841	A
1	1	1842	A
1	1	1849	C
1	1	1857	C
1	1	1858	A
1	1	1859	A
1	1	1863	G
1	1	1866	C
1	1	1878	G

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Mol	Chain	Res	Type
1	1	1879	A
1	1	1880	U
1	1	1906	G
1	1	1909	A
1	1	1937	U
1	1	1952	G
1	1	1953	G
1	1	2094	C
1	1	2100	A
1	1	2102	U
1	1	2112	U
1	1	2113	A
1	1	2114	C
1	1	2120	A
1	1	2121	G
1	1	2122	G
1	1	2131	A
1	1	2158	A
1	1	2163	C
1	1	2167	A
1	1	2168	A
1	1	2169	G
1	1	2191	U
1	1	2196	C
1	1	2197	C
1	1	2198	A
1	1	2205	U
1	1	2208	A
1	1	2209	U
1	1	2210	G
1	1	2223	A
1	1	2224	A
1	1	2242	A
1	1	2244	A
1	1	2247	G
1	1	2249	G
1	1	2250	G
1	1	2252	A
1	1	2255	A
1	1	2261	G
1	1	2262	A
1	1	2264	U

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Mol	Chain	Res	Type
1	1	2268	U
1	1	2269	U
1	1	2270	A
1	1	2272	G
1	1	2274	U
1	1	2277	C
1	1	2278	C
1	1	2316	G
1	1	2317	A
1	1	2318	U
1	1	2336	U
1	1	2363	A
1	1	2371	G
1	1	2378	C
1	1	2388	U
1	1	2393	G
1	1	2401	A
1	1	2402	A
1	1	2404	A
1	1	2414	G
1	1	2417	U
1	1	2418	G
1	1	2419	A
1	1	2433	U
1	1	2441	A
1	1	2442	G
1	1	2444	C
1	1	2502	A
1	1	2503	G
1	1	2514	U
1	1	2515	A
1	1	2522	G
1	1	2523	A
1	1	2526	C
1	1	2530	G
1	1	2531	C
1	1	2534	G
1	1	2538	U
1	1	2542	U
1	1	2543	U
1	1	2544	U
1	1	2545	C

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Mol	Chain	Res	Type
1	1	2546	C
1	1	2547	A
1	1	2549	G
1	1	2552	C
1	1	2555	G
1	1	2561	A
1	1	2562	A
1	1	2572	C
1	1	2573	G
1	1	2587	U
1	1	2593	A
1	1	2594	C
1	1	2606	G
1	1	2607	G
1	1	2614	G
1	1	2615	G
1	1	2621	G
1	1	2623	G
1	1	2625	C
1	1	2626	A
1	1	2635	A
1	1	2642	A
1	1	2651	G
1	1	2653	C
1	1	2654	C
1	1	2655	U
1	1	2656	A
1	1	2657	A
1	1	2659	G
1	1	2667	A
1	1	2674	A
1	1	2677	G
1	1	2688	U
1	1	2690	G
1	1	2693	C
1	1	2694	A
1	1	2696	A
1	1	2699	G
1	1	2702	A
1	1	2703	A
1	1	2704	A
1	1	2712	U

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Mol	Chain	Res	Type
1	1	2713	U
1	1	2714	G
1	1	2715	A
1	1	2718	U
1	1	2719	U
1	1	2721	A
1	1	2726	C
1	1	2727	A
1	1	2728	G
1	1	2729	U
1	1	2730	G
1	1	2732	G
1	1	2749	G
1	1	2752	U
1	1	2754	G
1	1	2758	A
1	1	2759	U
1	1	2762	A
1	1	2767	U
1	1	2770	G
1	1	2771	U
1	1	2772	C
1	1	2777	G
1	1	2778	G
1	1	2779	A
1	1	2790	A
1	1	2794	G
1	1	2795	U
1	1	2796	G
1	1	2798	C
1	1	2803	A
1	1	2810	C
1	1	2817	A
1	1	2818	U
1	1	2819	A
1	1	2820	A
1	1	2821	C
1	1	2822	U
1	1	2824	G
1	1	2825	C
1	1	2826	U
1	1	2838	A

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Mol	Chain	Res	Type
1	1	2842	U
1	1	2843	U
1	1	2857	C
1	1	2858	U
1	1	2861	U
1	1	2863	G
1	1	2865	U
1	1	2866	U
1	1	2867	C
1	1	2868	U
1	1	2869	U
1	1	2870	C
1	1	2872	A
1	1	2873	U
1	1	2875	U
1	1	2876	C
1	1	2877	G
1	1	2878	G
1	1	2879	C
1	1	2887	A
1	1	2889	C
1	1	2898	G
1	1	2901	G
1	1	2911	A
1	1	2924	U
1	1	2925	C
1	1	2926	A
1	1	2928	C
1	1	2930	A
1	1	2935	U
1	1	2936	A
1	1	2952	G
1	1	2953	U
1	1	2954	U
1	1	2955	U
1	1	2956	A
1	1	2965	U
1	1	2970	C
1	1	2971	A
1	1	2972	G
1	1	2980	U
1	1	2981	U

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Mol	Chain	Res	Type
1	1	2982	A
1	1	2983	C
1	1	2996	U
1	1	2997	G
1	1	3003	G
1	1	3012	A
1	1	3017	A
1	1	3021	A
1	1	3022	G
1	1	3023	U
1	1	3026	G
1	1	3028	G
1	1	3029	A
1	1	3030	G
1	1	3031	G
1	1	3032	A
1	1	3059	G
1	1	3078	U
1	1	3079	U
1	1	3086	A
1	1	3092	C
1	1	3093	C
1	1	3100	U
1	1	3101	G
1	1	3109	G
1	1	3129	A
1	1	3130	A
1	1	3131	U
1	1	3142	A
1	1	3143	C
1	1	3158	G
1	1	3165	A
1	1	3172	A
1	1	3173	G
1	1	3174	A
1	1	3176	G
1	1	3179	U
1	1	3181	C
1	1	3187	A
1	1	3196	U
1	1	3198	U
1	1	3207	U

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Mol	Chain	Res	Type
1	1	3213	A
1	1	3217	C
1	1	3218	A
1	1	3219	G
1	1	3229	G
1	1	3245	A
1	1	3247	G
1	1	3253	G
1	1	3259	U
1	1	3260	G
1	1	3270	U
1	1	3273	A
1	1	3276	G
1	1	3281	U
1	1	3287	U
1	1	3304	U
1	1	3316	A
1	1	3319	U
1	1	3334	U
1	1	3335	A
1	1	3341	U
1	1	3342	A
1	1	3343	G
1	1	3345	G
1	1	3350	C
1	1	3351	U
1	1	3352	U
1	1	3360	C
1	1	3368	U
1	1	3369	G
1	1	3375	A
1	1	3378	C
1	1	3389	U
2	2	16	G
2	2	23	U
2	2	34	U
2	2	35	C
2	2	39	G
2	2	59	A
2	2	62	C
2	2	63	G
2	2	72	A

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Mol	Chain	Res	Type
2	2	79	A
2	2	80	A
2	2	82	U
2	2	83	C
2	2	86	U
2	2	87	G
2	2	90	U
2	2	95	G
2	2	100	U
2	2	104	A
2	2	106	C
2	2	111	A
2	2	113	U
2	2	116	G
2	2	124	G
2	2	129	C
3	3	7	G
3	3	13	A
3	3	53	U
3	3	54	U
3	3	55	A
3	3	65	G
3	3	74	C
3	3	75	G
3	3	86	U
3	3	87	G
3	3	91	G
3	3	93	C
3	3	98	C
3	3	102	A
3	3	112	G
3	3	120	C
3	3	121	U

All (83) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	1	13	A
1	1	40	A
1	1	73	C
1	1	75	G
1	1	160	G

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Mol	Chain	Res	Type
1	1	239	G
1	1	284	A
1	1	547	G
1	1	619	A
1	1	645	A
1	1	649	A
1	1	720	A
1	1	765	C
1	1	849	C
1	1	916	G
1	1	975	C
1	1	1047	A
1	1	1064	A
1	1	1097	G
1	1	1102	A
1	1	1103	A
1	1	1128	U
1	1	1190	A
1	1	1205	A
1	1	1244	A
1	1	1302	A
1	1	1329	U
1	1	1355	A
1	1	1581	C
1	1	1607	U
1	1	1816	A
1	1	2101	C
1	1	2166	A
1	1	2204	C
1	1	2209	U
1	1	2260	U
1	1	2261	G
1	1	2263	C
1	1	2269	U
1	1	2271	A
1	1	2316	G
1	1	2317	A
1	1	2392	C
1	1	2501	U
1	1	2513	U
1	1	2525	G
1	1	2537	U

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Mol	Chain	Res	Type
1	1	2541	U
1	1	2554	A
1	1	2586	G
1	1	2593	A
1	1	2624	G
1	1	2625	C
1	1	2641	U
1	1	2652	U
1	1	2658	G
1	1	2725	U
1	1	2728	G
1	1	2758	A
1	1	2761	G
1	1	2770	G
1	1	2802	A
1	1	2857	C
1	1	2866	U
1	1	2868	U
1	1	2869	U
1	1	2875	U
1	1	2952	G
1	1	2954	U
1	1	2979	U
1	1	3030	G
1	1	3078	U
1	1	3157	U
1	1	3195	U
1	1	3218	A
1	1	3228	C
1	1	3269	U
1	1	3342	A
1	1	3350	C
1	1	3351	U
2	2	123	G
2	2	128	U
3	3	52	G

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

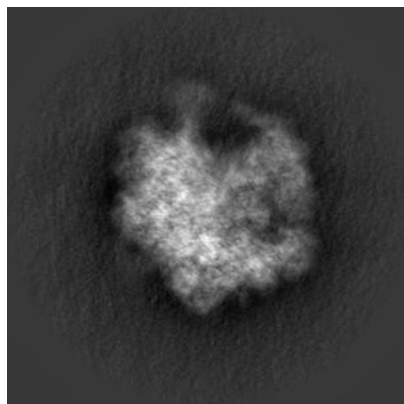
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-12892. These allow visual inspection of the internal detail of the map and identification of artifacts.

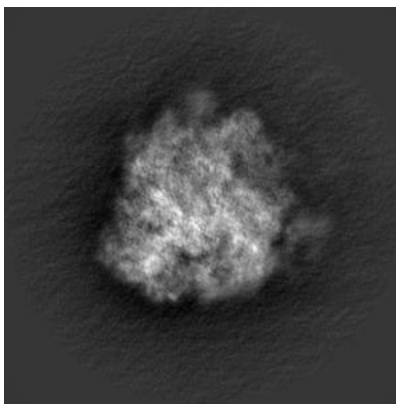
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

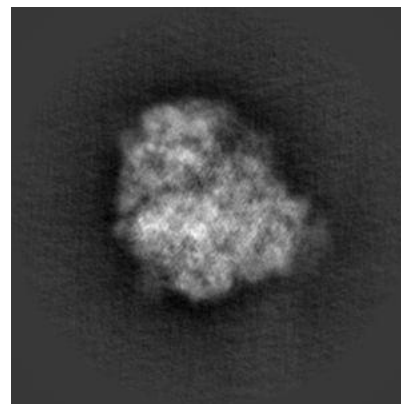
6.1.1 Primary map



X

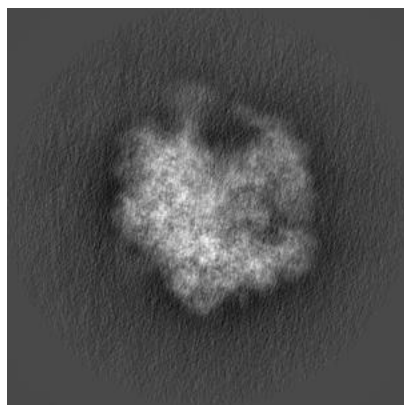


Y

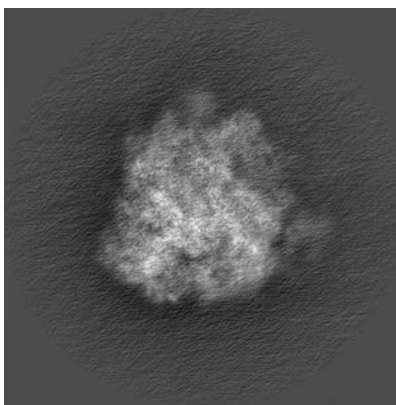


Z

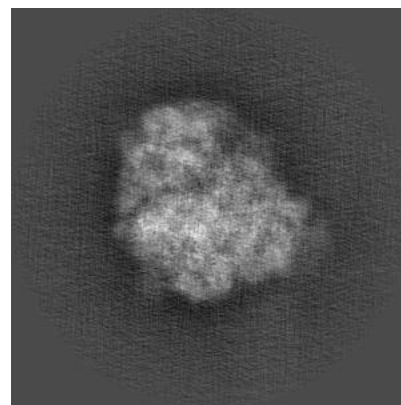
6.1.2 Raw map



X



Y

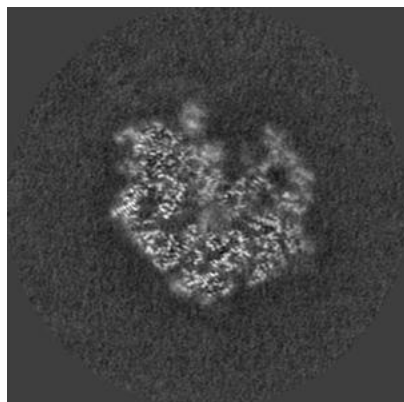


Z

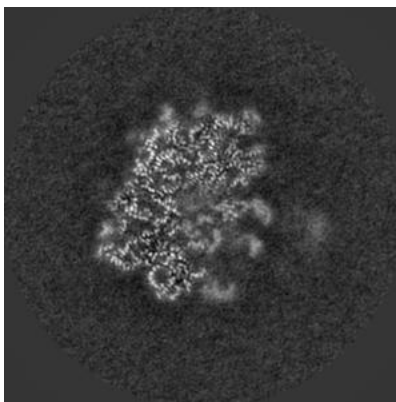
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

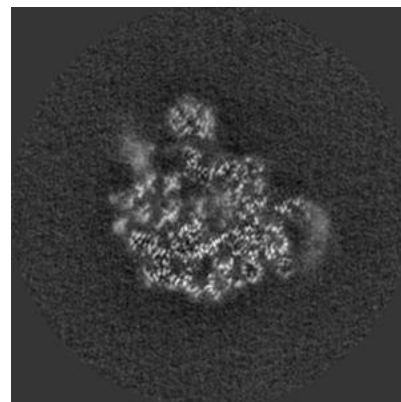
6.2.1 Primary map



X Index: 200

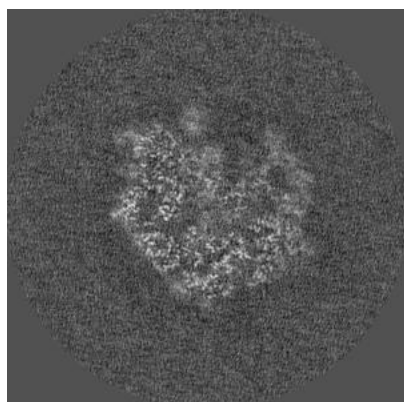


Y Index: 200

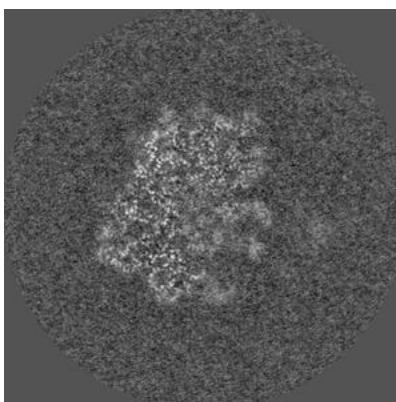


Z Index: 200

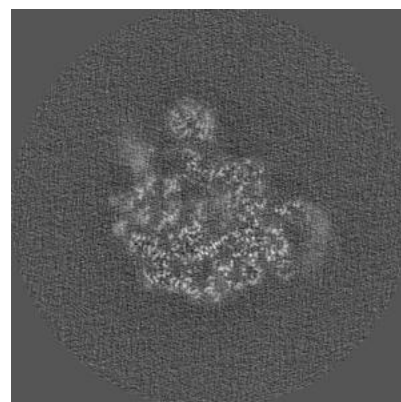
6.2.2 Raw map



X Index: 200



Y Index: 200

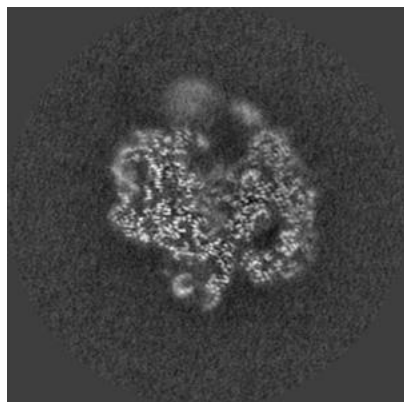


Z Index: 200

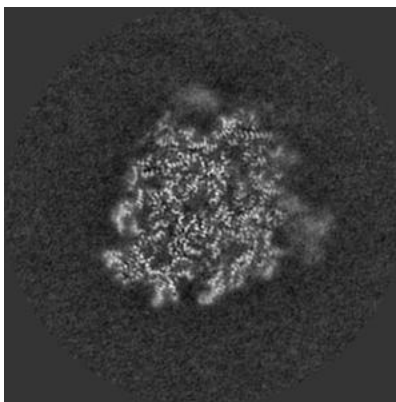
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

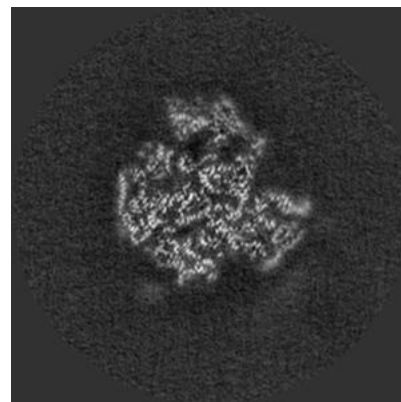
6.3.1 Primary map



X Index: 184

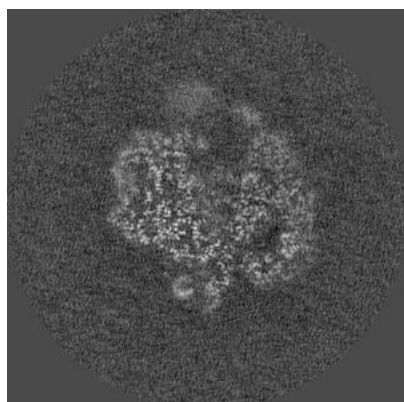


Y Index: 177

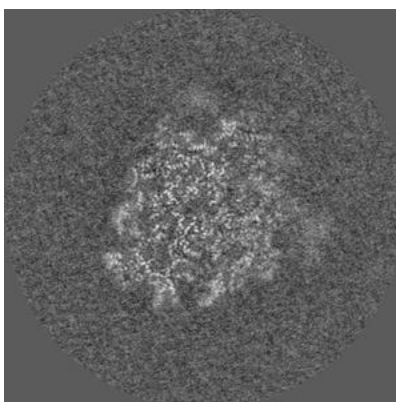


Z Index: 166

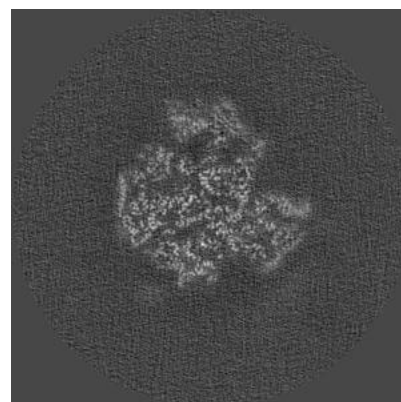
6.3.2 Raw map



X Index: 184



Y Index: 177

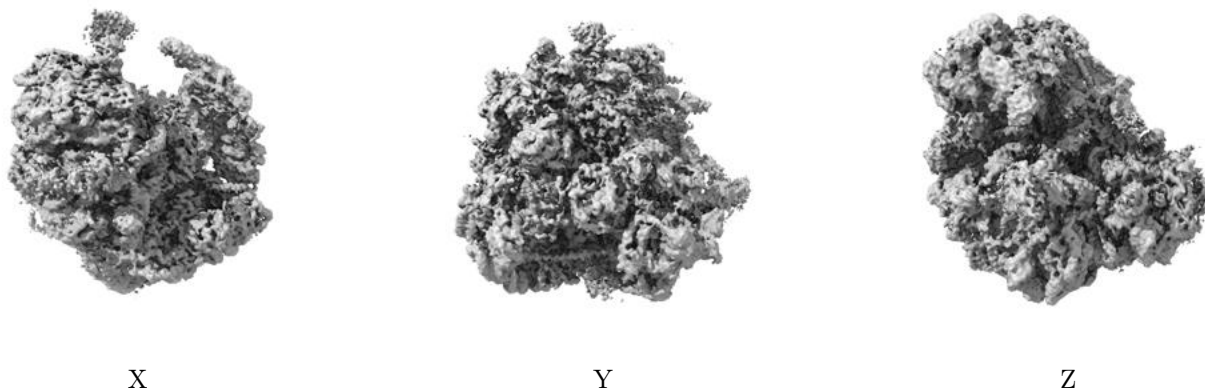


Z Index: 166

The images above show the largest variance slices of the map in three orthogonal directions.

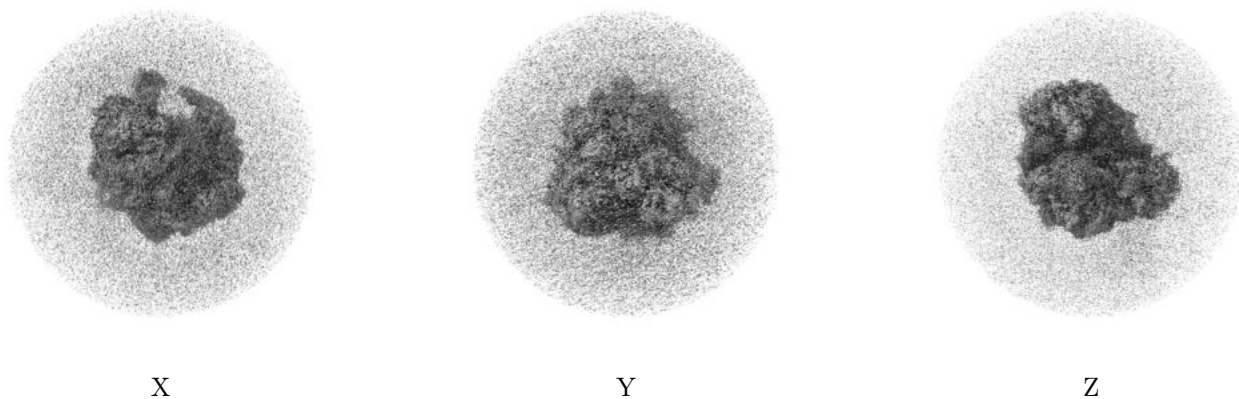
6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.026. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

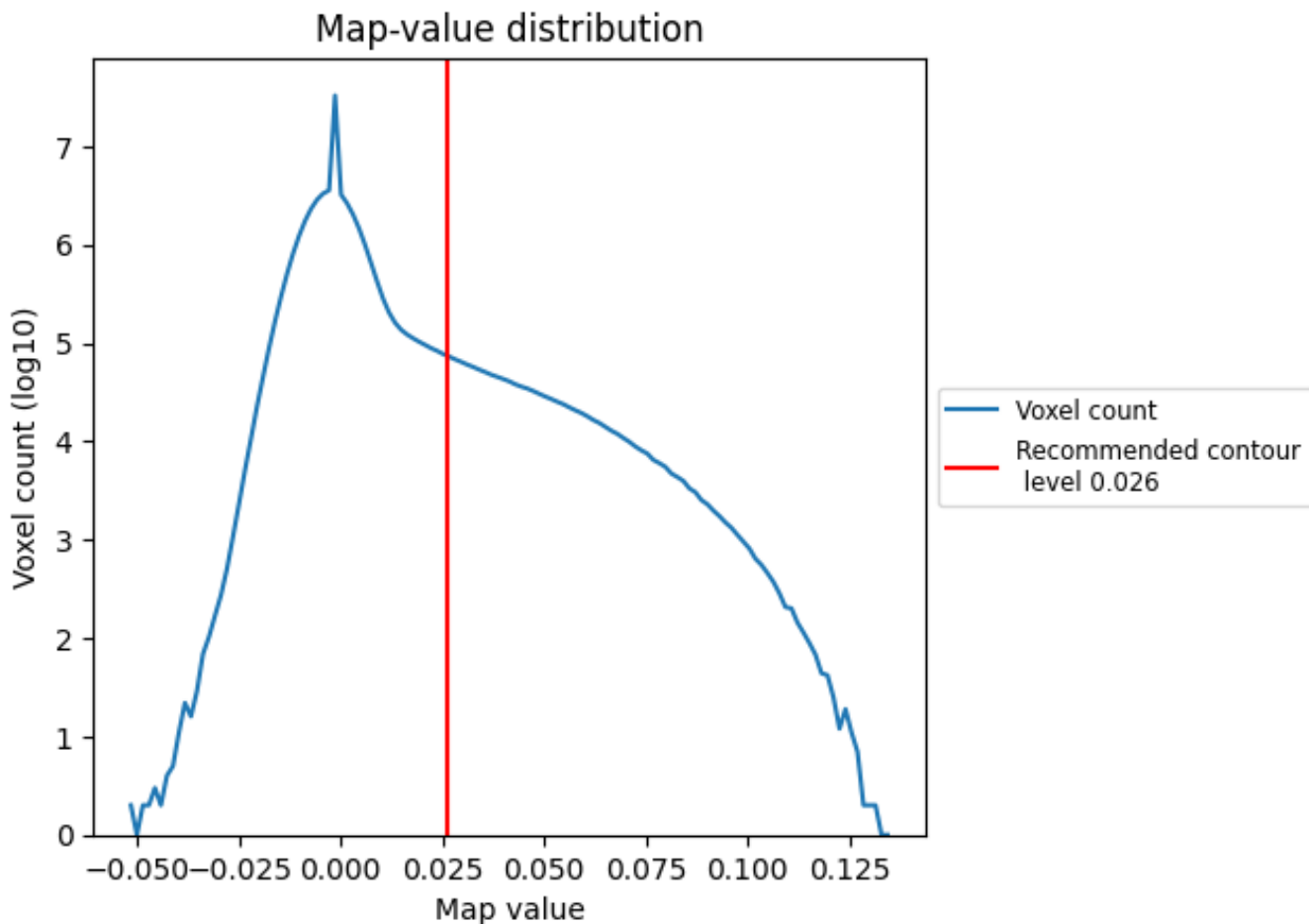
6.5 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

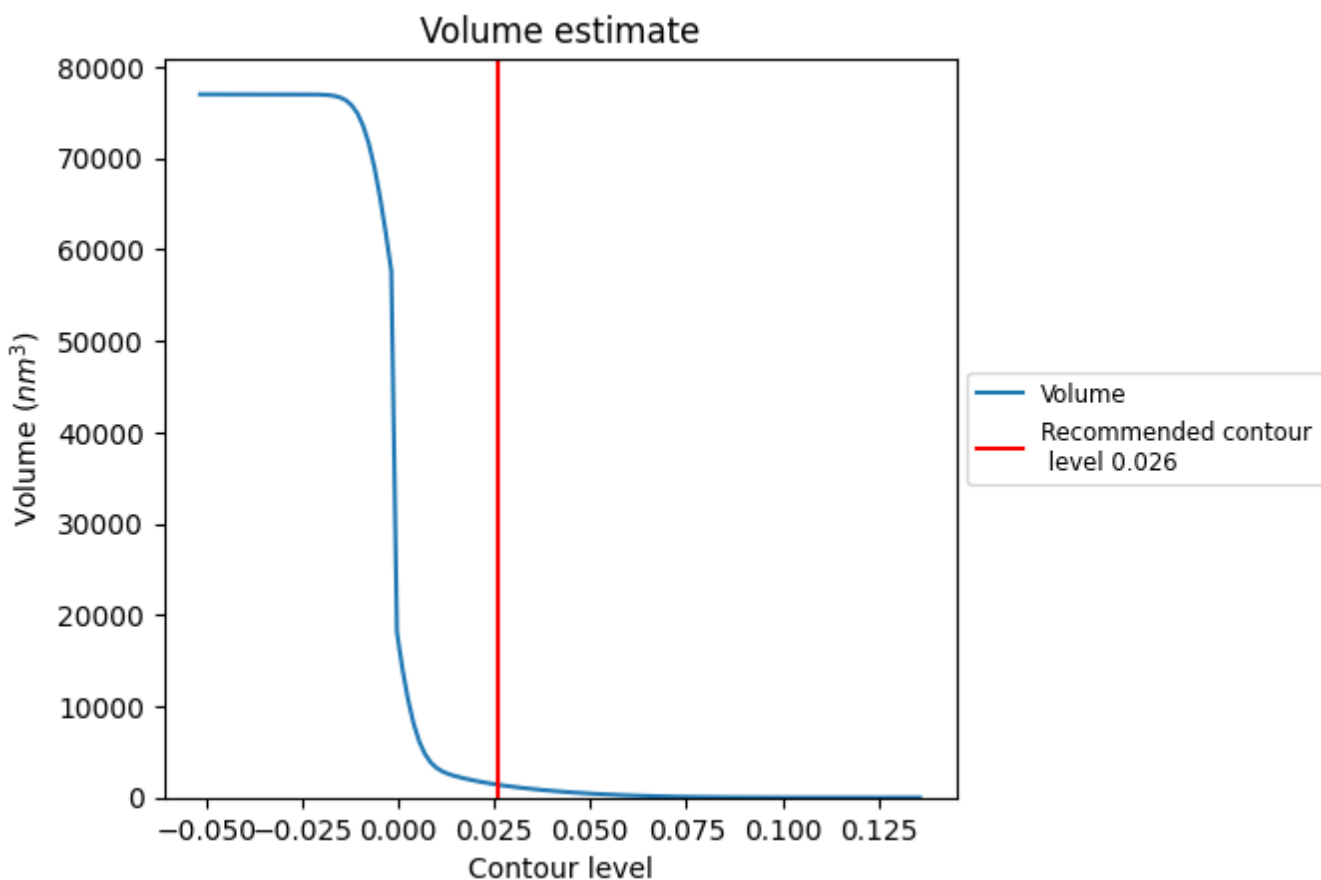
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

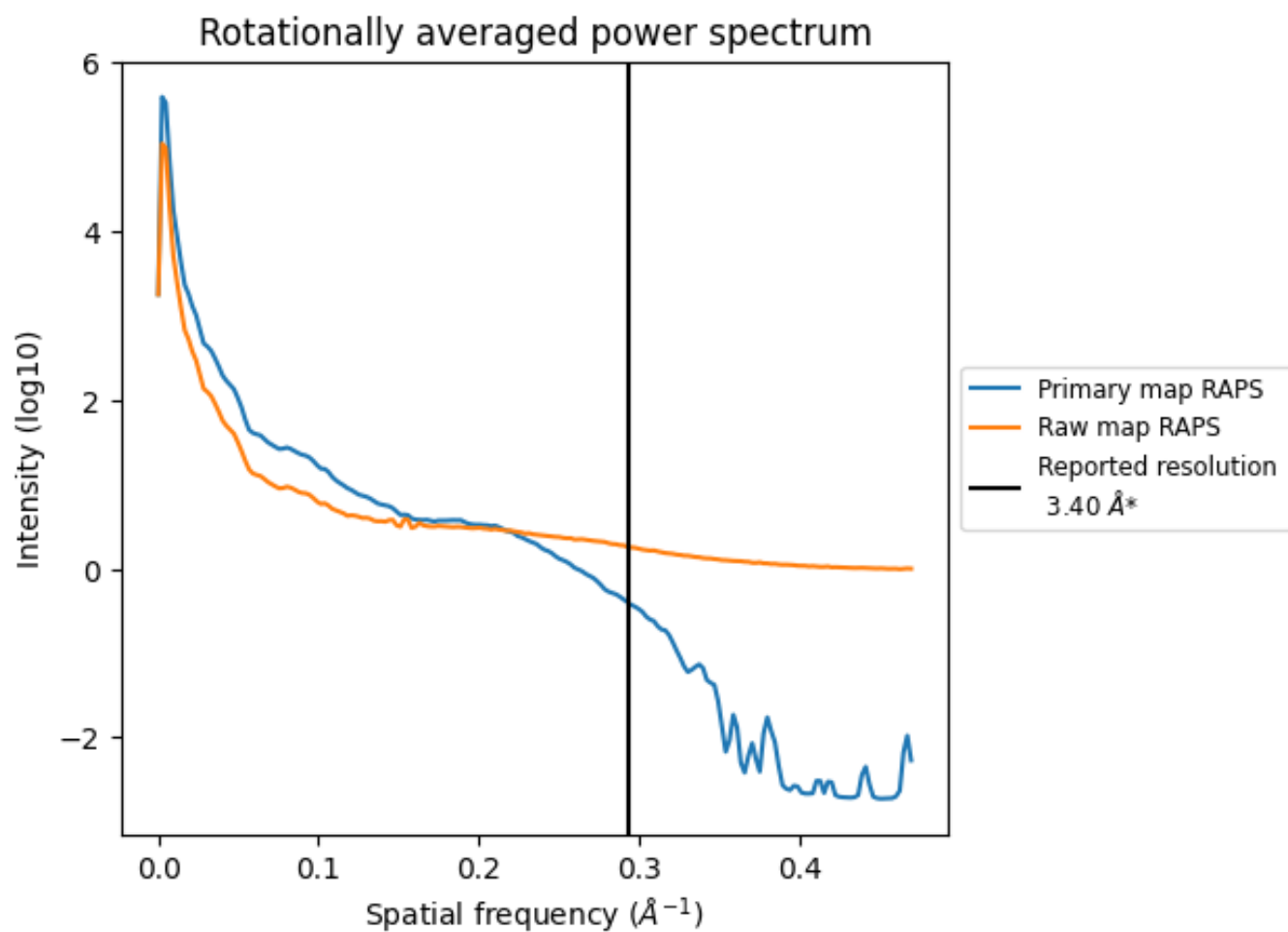
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1405 nm³; this corresponds to an approximate mass of 1269 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum i

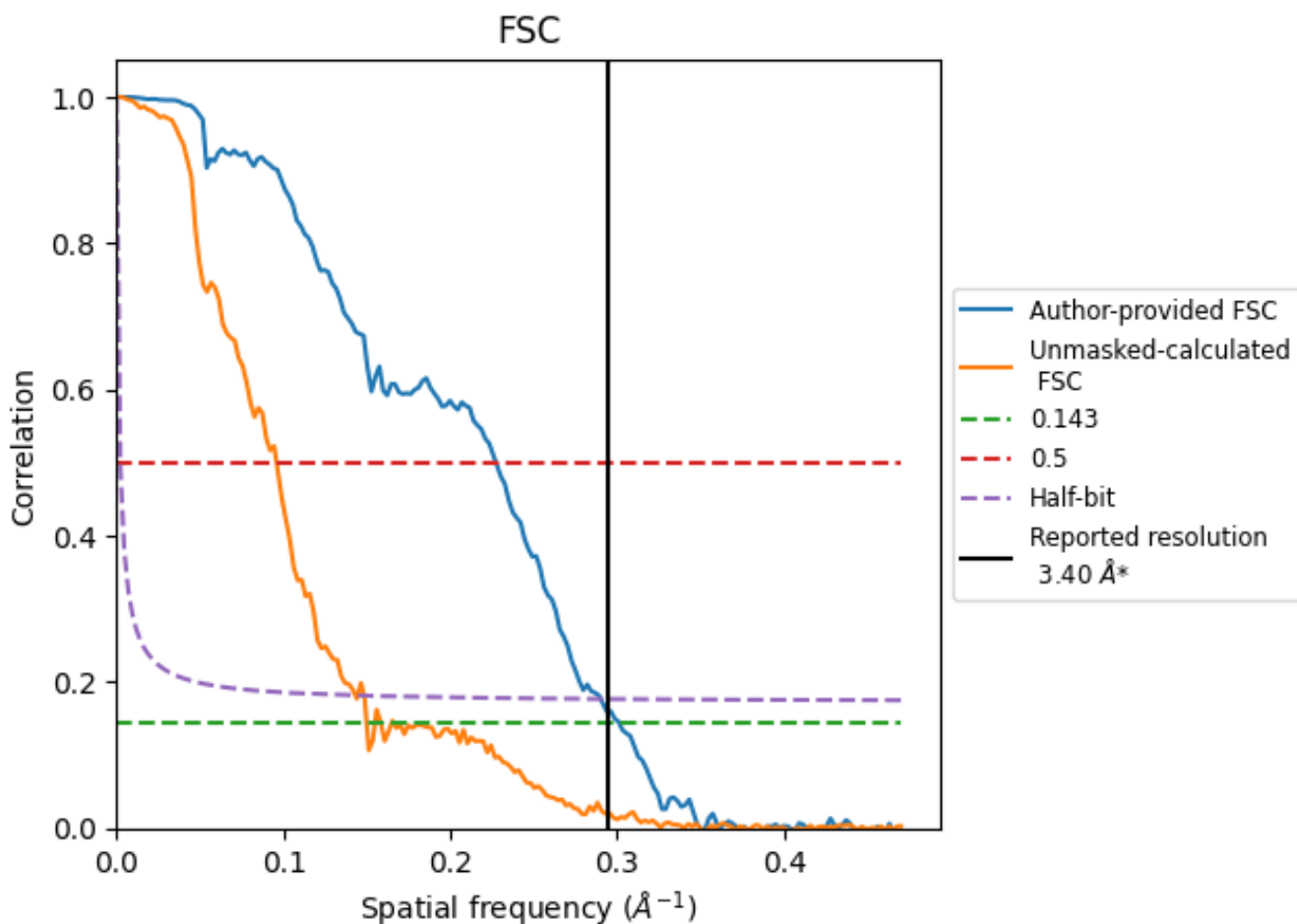


*Reported resolution corresponds to spatial frequency of 0.294 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.294 Å⁻¹

8.2 Resolution estimates [i](#)

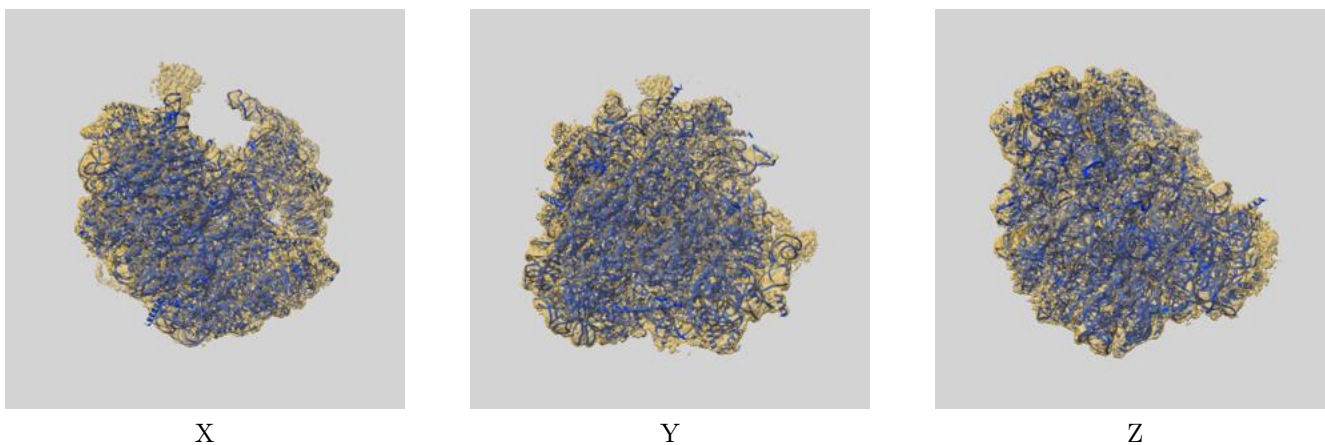
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.40	-	-
Author-provided FSC curve	3.32	4.40	3.45
Unmasked-calculated*	6.67	10.43	6.98

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 6.67 differs from the reported value 3.4 by more than 10 %

9 Map-model fit [i](#)

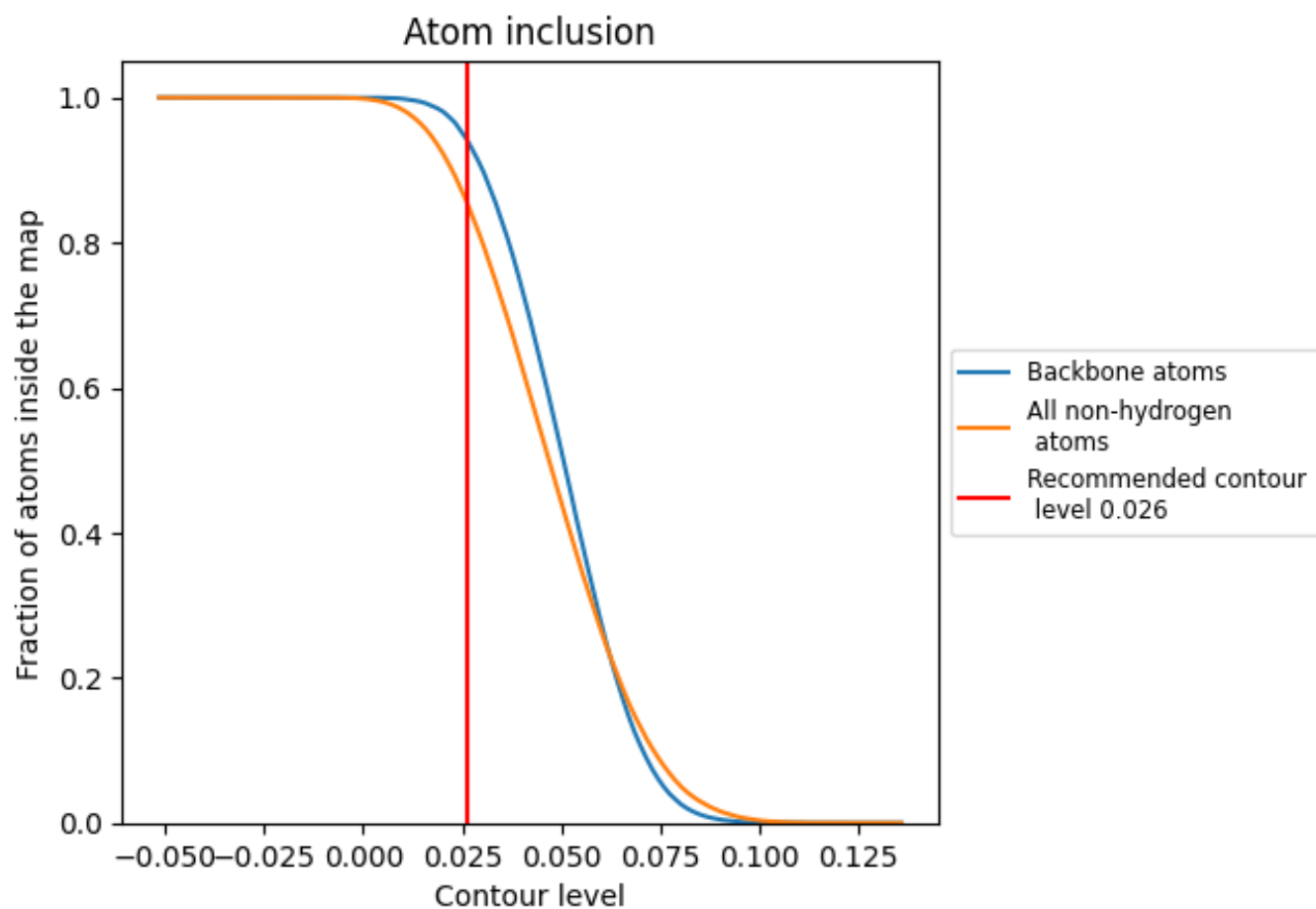
This section contains information regarding the fit between EMDB map EMD-12892 and PDB model 7OH3. Per-residue inclusion information can be found in section 3 on page 13.

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.026 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 86% of all non-hydrogen atoms, are inside the map.